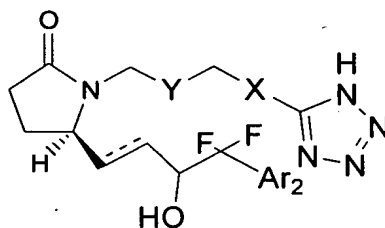


Case MC058YP

In the Claims

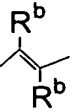
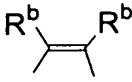
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17. A compound of structural formula I:



I

or a pharmaceutically acceptable salt, enantiomer, diastereomer, pro drug or mixture thereof, wherein

X is $(\text{CH}_2)_n$, O or S;

Y represents $(\text{C}(\text{R}^b)_2)_n$, triple bond,  or  ;

Case MC058YP

Ar₂ independently represent (CH₂)_mC₆₋₁₀aryl, (CH₂)_mC₅₋₁₀heteroaryl, (CH₂)_mC₃₋₁₀heterocycloalkyl, (CH₂)_mC₃₋₈cycloalkyl said cycloalkyl, heterocycloalkyl, aryl or heteroaryl unsubstituted or substituted with 1-3 groups of R_a;

R_a represents C₁₋₆alkoxy, C₁₋₆alkyl, CF₃, nitro, amino, cyano, C₁₋₆alkylamino, or halogen;

R^b independently represents H, halogen, C₁₋₆alkyl, C₃₋₆cylcoalkyl or

--- represents a double or single bond;

n represents 0-4; and

m represents 0-8.

18. The compound according to claim 17 wherein X and Y are (CH₂)_n, --- represents a double bond; and Ar₂ is phenyl.

19. The compound according to claim 18 wherein X is (CH₂)_n and n is 1 and Y is (CH₂)_n and n is 3.